



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 08:34 pm BST

PDB ID : 5BTE
Title : Crystal structure of Ashbya gossypii Rai1 in complex with pU(S)6-Mn2+
Authors : Wang, V.Y.; Tong, L.
Deposited on : 2015-06-03
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

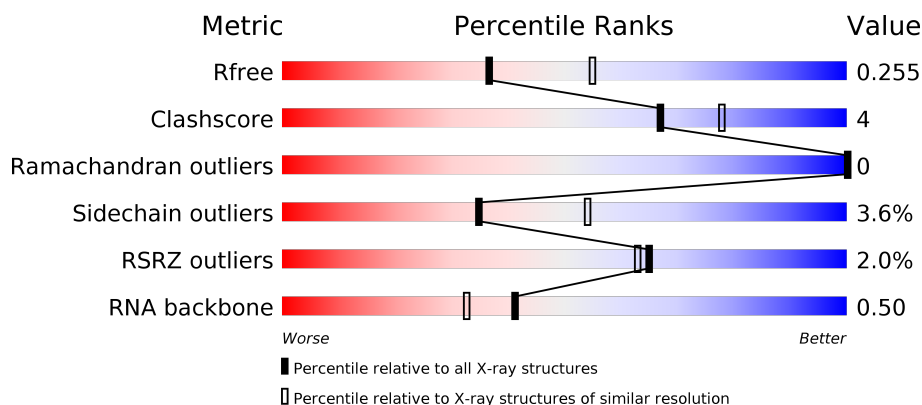
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)
RNA backbone	3102	1174 (2.80-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	376	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>10%</div> <div>..</div> </div> </div>
1	B	376	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div>..</div> </div> </div>
2	C	6	<div> <div>50%</div> <div>50%</div> </div>
2	D	6	<div> <div>50%</div> <div>50%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6550 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called AFR263Cp.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	368	Total	C	N	O	S	0	0	0
			2981	1901	520	550	10			
1	A	365	Total	C	N	O	S	0	0	0
			2954	1882	517	545	10			

- Molecule 2 is a RNA chain called RNA (5'-R(P*UP*(U37)P*(U37)P*UP*UP*U)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	6	Total	C	N	O	P	S	0	0	0
			121	54	12	47	6	2			
2	C	6	Total	C	N	O	P	S	0	0	0
			121	54	12	47	6	2			

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Mn	0	0
			3	3		
3	A	1	Total	Mn	0	0
			1	1		
3	C	2	Total	Mn	0	0
			2	2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

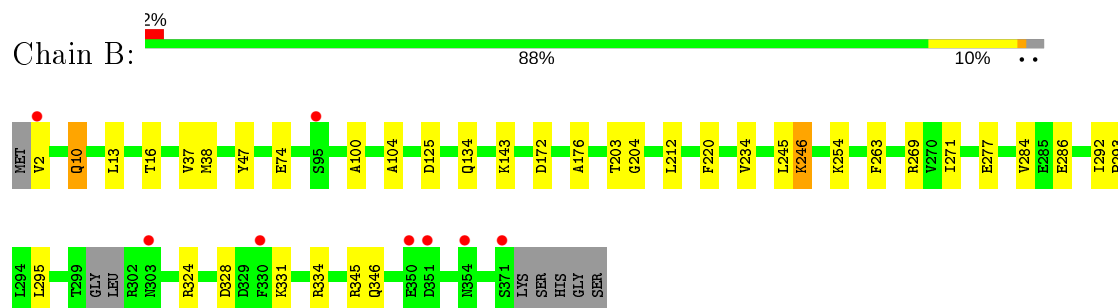
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	162	Total	O	0	0
			162	162		
5	D	16	Total	O	0	0
			16	16		
5	C	14	Total	O	0	0
			14	14		
5	A	165	Total	O	0	0
			165	165		

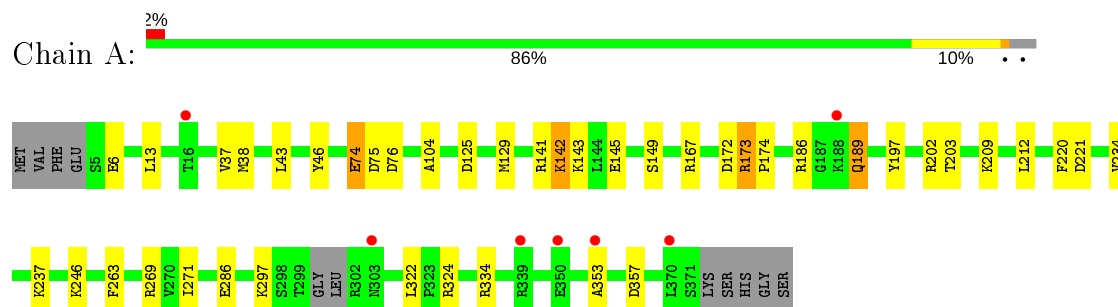
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: AFR263Cp



- Molecule 1: AFR263Cp



- Molecule 2: RNA (5'-R(P*UP*(U37)P*(U37)P*UP*UP*U)-3')



- Molecule 2: RNA (5'-R(P*UP*(U37)P*(U37)P*UP*UP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.66 Å 79.74 Å 160.14 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.41 – 2.40 42.37 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.41-2.40) 99.7 (42.37-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.62 (at 2.39 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.196 , 0.254 0.203 , 0.255	Depositor DCC
R_{free} test set	1787 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6550	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 51.78 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.3413e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: U37, MN, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.66	0/3016	0.80	1/4065 (0.0%)
1	B	0.63	0/3044	0.77	1/4103 (0.0%)
2	C	1.36	2/87 (2.3%)	0.76	0/128
2	D	1.19	1/87 (1.1%)	0.95	1/128 (0.8%)
All	All	0.67	3/6234 (0.0%)	0.79	3/8424 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	U	OP3-P	-10.37	1.48	1.61
2	C	2	U	OP3-P	-9.58	1.49	1.61
2	C	5	U	O3'-P	-7.54	1.52	1.61

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	U	O5'-P-OP2	-6.85	99.54	105.70
1	A	167	ARG	NE-CZ-NH2	6.69	123.64	120.30
1	B	324	ARG	NE-CZ-NH1	5.68	123.14	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2954	0	2946	24	0
1	B	2981	0	2970	23	0
2	C	121	0	61	1	0
2	D	121	0	61	2	0
3	A	1	0	0	0	0
3	B	3	0	0	0	0
3	C	2	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	165	0	0	4	1
5	B	162	0	0	6	2
5	C	14	0	0	1	0
5	D	16	0	0	1	1
All	All	6550	0	6038	48	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (48) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ASP:OD1	1:A:334:ARG:NH1	2.06	0.89
1:B:176:ALA:O	5:B:501:HOH:O	2.04	0.75
2:D:3:U37:S	5:D:114:HOH:O	2.45	0.74
1:A:172:ASP:OD2	5:A:501:HOH:O	2.09	0.70
1:B:10:GLN:OE1	5:B:502:HOH:O	2.12	0.67
1:B:234:VAL:HG21	1:B:271:ILE:HD12	1.77	0.67
1:B:16:THR:O	1:B:16:THR:HG22	1.97	0.65
1:B:172:ASP:OD2	5:B:503:HOH:O	2.15	0.64
1:B:220:PHE:CD2	1:B:269:ARG:HD3	2.37	0.59
1:A:234:VAL:HG21	1:A:271:ILE:HD12	1.86	0.57
1:B:125:ASP:OD1	1:B:334:ARG:NH1	2.39	0.56
1:A:186:ARG:HA	1:A:189:GLN:HE21	1.75	0.52
1:A:37:VAL:HG12	1:A:38:MET:HG2	1.93	0.51
1:A:322:LEU:O	1:A:324:ARG:HG3	2.12	0.50
1:A:186:ARG:HA	1:A:189:GLN:HG2	1.94	0.50
1:B:292:ILE:N	1:B:293:PRO:CD	2.75	0.49
1:A:141:ARG:NH1	5:A:514:HOH:O	2.44	0.49
1:B:328:ASP:OD1	1:B:331:LYS:NZ	2.46	0.48
1:A:203:THR:HG21	5:A:557:HOH:O	2.14	0.47
1:B:245:LEU:HB3	1:B:246:LYS:HE3	1.96	0.47
1:B:104:ALA:O	1:B:212:LEU:HA	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ALA:O	1:B:134:GLN:HG2	2.17	0.45
1:A:220:PHE:HB2	1:A:234:VAL:HB	1.98	0.45
1:B:47:TYR:HE1	1:A:74:GLU:HG2	1.81	0.45
1:A:173:ARG:HB2	1:A:174:PRO:CD	2.46	0.44
1:B:331:LYS:HD3	1:B:346:GLN:HE22	1.82	0.44
1:B:16:THR:O	1:B:16:THR:CG2	2.65	0.44
1:A:221:ASP:OD1	1:A:221:ASP:N	2.49	0.44
1:A:173:ARG:HB2	1:A:174:PRO:HD2	2.00	0.43
1:A:220:PHE:CD2	1:A:269:ARG:CD	3.01	0.43
1:B:271:ILE:HG23	1:B:284:VAL:HG13	2.01	0.43
1:B:203:THR:HG21	5:B:560:HOH:O	2.18	0.43
1:B:203:THR:OG1	1:B:204:GLY:N	2.51	0.43
1:A:129:MET:HE1	1:A:353:ALA:HA	2.01	0.42
1:B:37:VAL:HG12	1:B:38:MET:HG2	2.02	0.42
1:A:76:ASP:HA	1:A:142:LYS:O	2.19	0.42
2:C:4:U37:O5'	2:C:4:U37:H6	2.20	0.42
1:B:254:LYS:HE2	2:D:4:U37:S	2.59	0.42
1:A:104:ALA:O	1:A:212:LEU:HA	2.19	0.42
1:A:43:LEU:HD13	1:A:197:TYR:CE2	2.55	0.41
1:B:220:PHE:HB2	1:B:234:VAL:HB	2.02	0.41
1:A:46:TYR:CZ	1:A:271:ILE:HD13	2.55	0.41
5:B:599:HOH:O	1:A:75:ASP:HA	2.21	0.41
1:B:220:PHE:CD2	1:B:269:ARG:CD	3.03	0.40
1:A:357:ASP:HB2	5:A:596:HOH:O	2.21	0.40
1:A:202:ARG:NH2	1:A:209:LYS:HD3	2.36	0.40
5:C:508:HOH:O	1:A:237:LYS:HA	2.22	0.40
1:B:277:GLU:HG2	5:B:509:HOH:O	2.20	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:643:HOH:O	5:A:658:HOH:O[1_455]	1.95	0.25
5:B:501:HOH:O	5:D:108:HOH:O[3_454]	2.17	0.03

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	361/376 (96%)	350 (97%)	11 (3%)	0	100	100
1	B	364/376 (97%)	351 (96%)	13 (4%)	0	100	100
All	All	725/752 (96%)	701 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	315/324 (97%)	302 (96%)	13 (4%)	30	48
1	B	318/324 (98%)	308 (97%)	10 (3%)	40	60
All	All	633/648 (98%)	610 (96%)	23 (4%)	35	54

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2	VAL
1	B	10	GLN
1	B	13	LEU
1	B	74	GLU
1	B	143	LYS
1	B	246	LYS
1	B	263	PHE
1	B	286	GLU

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Mol	Chain	Res	Type
1	B	295	LEU
1	B	345	ARG
1	A	6	GLU
1	A	13	LEU
1	A	74	GLU
1	A	142	LYS
1	A	143	LYS
1	A	145	GLU
1	A	149	SER
1	A	173	ARG
1	A	189	GLN
1	A	246	LYS
1	A	263	PHE
1	A	286	GLU
1	A	297	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	108	GLN
1	B	343	HIS
1	B	346	GLN
1	A	108	GLN
1	A	189	GLN
1	A	253	GLN
1	A	343	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	3/6 (50%)	0	0
2	D	3/6 (50%)	0	0
All	All	6/12 (50%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	U37	C	3	2	14,21,22	1.12	2 (14%)	13,30,33	0.69	0
2	U37	D	4	2	14,21,22	1.32	3 (21%)	13,30,33	0.76	0
2	U37	D	3	2	14,21,22	1.09	0	13,30,33	0.85	0
2	U37	C	4	2	14,21,22	1.06	1 (7%)	13,30,33	0.43	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	U37	C	3	2	-	0/5/25/26	0/2/2/2
2	U37	D	4	2	-	0/5/25/26	0/2/2/2
2	U37	D	3	2	-	0/5/25/26	0/2/2/2
2	U37	C	4	2	-	0/5/25/26	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	4	U37	C2-N3	-2.58	1.33	1.38
2	C	4	U37	C2-N3	-2.26	1.33	1.38
2	D	4	U37	C2'-C1'	-2.21	1.50	1.53
2	C	3	U37	C2-N3	-2.19	1.33	1.38
2	D	4	U37	O4'-C1'	2.14	1.44	1.41
2	C	3	U37	C6-N1	-2.13	1.33	1.35

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	4	U37	1	0
2	D	3	U37	1	0
2	C	4	U37	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	404	-	4,4,4	0.26	0	6,6,6	0.63	0
4	SO4	A	402	-	4,4,4	0.53	0	6,6,6	0.93	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	365/376 (97%)	-0.37	7 (1%) 66 64	10, 24, 51, 71	0
1	B	368/376 (97%)	-0.37	8 (2%) 62 60	10, 25, 54, 77	0
2	C	4/6 (66%)	-1.18	0 100 100	17, 23, 24, 30	0
2	D	4/6 (66%)	-1.22	0 100 100	16, 20, 24, 32	0
All	All	741/764 (96%)	-0.38	15 (2%) 65 63	10, 24, 52, 77	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	VAL	3.6
1	B	95	SER	3.5
1	B	354	ASN	3.4
1	A	370	LEU	3.2
1	A	16	THR	3.1
1	A	339	ARG	2.9
1	A	353	ALA	2.7
1	B	303	ASN	2.6
1	B	351	ASP	2.4
1	A	350	GLU	2.3
1	A	303	ASN	2.2
1	B	350	GLU	2.2
1	B	371	SER	2.1
1	A	188	LYS	2.1
1	B	330	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	U37	C	3	20/21	0.95	0.11	13,15,17,28	0
2	U37	D	3	20/21	0.96	0.10	12,16,18,22	0
2	U37	C	4	20/21	0.97	0.10	17,19,21,21	0
2	U37	D	4	20/21	0.98	0.11	18,22,25,26	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MN	B	403	1/1	0.98	0.04	42,42,42,42	0
4	SO4	A	402	5/5	0.99	0.12	18,21,21,24	0
3	MN	A	401	1/1	0.99	0.02	43,43,43,43	0
4	SO4	B	404	5/5	0.99	0.13	20,22,22,23	0
3	MN	B	401	1/1	1.00	0.07	8,8,8,8	0
3	MN	C	401	1/1	1.00	0.05	10,10,10,10	0
3	MN	B	402	1/1	1.00	0.07	11,11,11,11	0
3	MN	C	402	1/1	1.00	0.09	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.